## AMENDMENTS TO THE CLAIMS

The following listing of claims replaces all prior listings of claims presented in the application.

## 1. (Currently amended) A compound of formula (I),

$$R_{1}$$
 $R_{1}$ 
 $R_{1}$ 
 $R_{1}$ 
 $R_{11}$ 
 $R_{12}$ 
 $R_{13}$ 
 $R_{14}$ 
 $R_{11}$ 
 $R_{12}$ 
 $R_{2}$ 
 $R_{3}$ 
 $R_{4}$ 
 $R_{5}$ 
 $R_{5}$ 
 $R_{6}$ 

and its tautomeric forms, its stereoisomers, and its pharmaceutically acceptable salts,

wherein R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, R<sub>8</sub>, R<sub>9</sub>, R<sub>10</sub>, R<sub>11</sub> and R<sub>12</sub> may be the same or different and each independently represent hydrogen, halogen, perhaloalkyl, hydroxy, amino, nitro, cyano, formyl, amidino, guanidino, substituted or unsubstituted groups such as selected from the group consisting of linear or branched (C<sub>1</sub>-C<sub>12</sub>)alkyl, (C<sub>2</sub>-C<sub>12</sub>)alkenyl, (C<sub>2</sub>-C<sub>12</sub>)alkynyl, (C<sub>3</sub>-C<sub>7</sub>)cycloalkyl, bicycloalkyl, bicycloalkenyl, (C<sub>1</sub>-C<sub>12</sub>)alkoxy, cyclo(C<sub>3</sub>-C<sub>7</sub>)alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heterocyclylalkyl, heteroaryloxy, heteroaryloxy, heterocyclylalkyloxy, acyl, acyloxy, acylamino, monoalkylamino, dialkylamino, aralkylamino, aralkylamino, alkoxycarbonyl, aryloxycarbonyl, aralkoxycarbonyl, heteroaryloxycarbonyl, heterocyclylalkoxycarbonyl, heteroaryloxycarbonyl, aminoalkyl,

monoalkylaminoalkyl, dialkylaminoalkyl, alkoxyalkyl, aryloxyalkyl, aralkoxyalkyl, alkylthio, aralkyloxycarbonylamino, alkoxycarbonylamino, aryloxycarbonylamino. thioalkyl, aminocarbonylamino, alkylaminocarbonylamino, dialkylaminocarbonylamino, alkylamidino, alkylguanidino, dialkylguanidino, hydrazino, and hydroxylamino; or the adjacent groups like R1 and R2 or R2 and R3 or R3 and R4 or R5 and R6 or R6 and R7 or R7 and R8 together with carbon atoms to which they are attached may form a 5, 6, or 7 membered ring, which may further optionally contain one or more double bonds and/or one or more heteroatoms such as selected from the group consisting of the group "Oxygen", "Nitrogen", "Sulfur"-or and "Selenium"-and combinations of double bond and heteroatoms; or R<sub>9</sub> and R<sub>10</sub> or R<sub>11</sub> and R<sub>12</sub> together represent double bond attached to "Oxygen" or "Sulfur"; or Ro and R10 or R11 and R12 together with the carbon atoms to which they are attached may form a 3, 4, 5, or 6 membered ring, which may further optionally contain one or more double bonds, and/or one or more heteroatoms such as selected from the group consisting of the group "Oxygen", "Nitrogen", "Sulfur" or and "Selenium," as above defined;

 $R_{13}$  and  $R_{14}$  may be the same or different and each independently represents hydrogen, substituted or unsubstituted groups such as selected from the group consisting of linear or branched  $(C_1-C_{12})$ alkyl,  $(C_2-C_{12})$ alkenyl,  $(C_2-C_{12})$ alkynyl,  $(C_2-C_{12})$ alkanoyl  $(C_3-C_7)$ cycloalkyl,  $(C_3-C_7)$ cycloalkenyl, bicycloalkenyl, aryl, aralkyl, heteroaryl,  $\theta$  and heterocyclylalkyl; or  $R_{13}$  and  $R_{14}$  along with the nitrogen atom, may form a 3, 4, 5, 6 or 7-membered heterocyclic ring, wherein the ring may be further substituted, and it may have either one, two or three double bonds or "additional heteroatoms"—as-defined-above; and

"n" is an integer ranging from 1 to 8, .

(Currently amended) A compound according to Claim 1, which is selected from the group consisting of:

6-(2-N,N-Dimethylaminoethyl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;

4-Bromo-6-(2-N,N-dimethylaminoethyl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;

4-Chloro-6-(2-N,N-dimethylaminoethyl)-benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;

 $6\hbox{-}(2\hbox{-}N,N\hbox{-}Dimethylaminoethyl)\hbox{-}4\hbox{-}fluor obenzo[d] is othiazolo[3,2\hbox{-}a] indol\hbox{-}S,S\hbox{-}dioxide;$ 

6-(2-N,N-Dimethylaminoethyl)-4-methylbenzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;

 $\label{eq:condition} 6-(2-N,N-Dimethylaminoethyl)-4-methylbenzo[d] is othiazolo[3,2-a] indol-S,S-dioxide hydrochloride salt;$ 

 $\label{eq:condition} 6-(2-N,N-Dimethylaminoethyl)-4-methylbenzo[d] is othiazolo[3,2-a] indol-S,S-dioxide maleate$ 

salt;

 $\label{eq:condition} 6-(2-N,N-Dimethylaminoethyl)-4-methylbenzo[d] isothiazolo[3,2-a] indol-S,S-dioxide D.L-malic acid salt:$ 

 $\label{eq:condition} 6-(2-N,N-Dimethylaminoethyl)-4-methylbenzo[d] is othiazolo[3,2-a] indol-S,S-dioxide oxalate salt:$ 

 $\label{eq:condition} 6-(2-N,N-Dimethylaminoethyl)-4-methylbenzo[d] isothiazolo[3,2-a] indol-S,S-dioxide citrate salt;$ 

 $6\hbox{-}(2\hbox{-}N,N\hbox{-}Dimethylaminoethyl)\hbox{-}4\hbox{-}methoxybenzo[d] isothiazolo[3,2\hbox{-}a] indol\hbox{-}S,S\hbox{-}dioxide;$ 

 $6\hbox{-}(2\hbox{-}N,N\hbox{-}Dimethylaminoethyl)\hbox{-}8\hbox{-}methoxybenzo[d] is othiazolo[3,2\hbox{-}a] indol\hbox{-}S,S\hbox{-}dioxide;$ 

4-Bromo-6-(2-N,N-dimethylaminoethyl)-8-methoxybenzo[d]isothiazolo[3,2-a]indol-S,S-

dioxide;

 $\label{lem:condition} 4- Chloro-6-(2-N,N-dimethylaminoethyl)-8-methoxybenzo[d] is othiazolo[3,2-a] indol-S, S-methoxybenzo[d] indol-S, S-methoxybenzo[d] is othiazolo[3,2-a] indol-S, S-methoxybenzo[d] indol-S$ 

dioxide;

 $\label{eq:condition} 6-(2-N,N-Dimethylaminoethyl)-4-fluoro-8-methoxybenzo[d] isothiazolo[3,2-a] indol-S, S-dioxide:$ 

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 $6\hbox{-}(2\hbox{-}N,N\hbox{-}Dimethylaminoethyl)-4-methyl-8-methoxybenzo[d] is othiazolo[3,2\hbox{-}a] indol-S,S\hbox{-}dioxide;$ 

- $\label{eq:condition} 6\mbox{-}(2\mbox{-N,N-Dimethylaminoethyl)-4,8-dimethoxybenzo[d]} is othiazolo[3,2-a] indol-S, S-dioxide;$ 
  - $6\hbox{-}(2\hbox{-}N,N\hbox{-}Dimethylaminoethyl)\hbox{-}2\hbox{-}ethylbenzo[d] is othiazolo[3,2\hbox{-}a] indol\hbox{-}S,S\hbox{-}dioxide;$
  - $\hbox{2-Chloro-6-(2-N,N-dimethylaminoethyl)} benzo[d] is othiazolo[3,2-a] indol-S,S-dioxide;$
- 2, 4- Dichloro-6-(2-N,N-dimethylaminoethyl)-benzo[d] is othiazolo[3,2-a] indol-S, S-dioxide;
- 2,3- Dichloro-6-(2-N,N-dimethylaminoethyl)-benzo[d] is othiazolo[3,2-a] indol-S,S-dimethylaminoethyl)-benzo[d] is othiazolo[3,2-a] indol-S,S-dimethylaminoethyl] is othiazolo[3,2-a] indol-S,S-dimethylaminoethyl] is othiazolo[3,2-a] indol-S,S-dimethylaminoethyl] is othiazolo[3,2-a] indol-S,S-dimethylaminoethylaminoethyl] is othiazolo[3,2-a] indol-S,S-dimethylaminoethy

dioxide;

5- Chloro-6-(2-N,N-dimethylaminoethyl)-2-methylbenzo[d] is othiazolo[3,2-a] indol-S, S-dimethylaminoethyl)-2-methylbenzo[d] is othiazolo[3,2-a] indol-S, S-dimethylaminoethyll-2-methylbenzo[d] is othiazolo[3,2-a] indol-S, S-dimethylaminoethyll-2-methylaminoethyll-2-methyll-2-methylaminoethyll-2-met

dioxide;

2,4,5-Trichloro-6-(2-N,N-dimethylaminoethyl)-benzo[d]isothiazolo[3,2-a]indol-S,S-

dioxide;

6-(2-N,N-Dimethylaminoethyl)-2,4-difluorobenzo[d]isothiazolo[3,2-a]indol-S,S-

dioxide;

- $\label{eq:condition} 6-(2-N,N-dimethylaminoethyl)-4-fluoro-8-methylbenzo[d] isothiazolo[3,2-a] indol-S, S-dioxide:$
- 2,4-Difluoro-6-(2-N,N-dimethylaminoethyl)-8-methylbenzo[d] isothiazolo[3,2-a] indol-S,S-dioxide;
  - $6\hbox{-}(2\hbox{-}N,N\hbox{-}Dimethylaminoethyl)\hbox{-}2\hbox{-}methoxybenzo[d] is othiazolo[3,2\hbox{-}a] indol\hbox{-}S,S\hbox{-}dioxide;$
- $\label{eq:continuous} 6\text{-}(2\text{-N,N-Dimethylaminoethyl})\text{-}2,\\ 8\text{-}dimethoxybenzo[d] is othiazolo[3,2\text{-}a] indol-S, S-dioxide:$

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  - 6-(2-N,N-Dimethylaminoethyl)-8-methylbenzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
- 6-(3-N,N-Dimethylamino-1-hydroxyprop-1-yl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide:
- $\label{lem:condition} 4-Bromo-6-(3-N,N-Dimethylamino-1-hydroxyprop-1-yl) benzo[d] is othiazolo [3,2-a] indol-S,S-dioxide;$
- $\label{lem:condition} 6-(3-N,N-Dimethylamino-1-hydroxyprop-1-yl)-8-methoxybenzo[d] isothiazolo[3,2-a] indol-S,S-dioxide;$
- $\label{eq:continuous} 6-(3-N,N-Dimethylamino-1-hydroxyprop-1-yl)-8-methylbenzo[d] is othiazolo [3,2-a] indol-S,S-dioxide;$
- 4-Bromo-6-(3-N,N-dimethylamino-1-hydroxyprop-1-yl)-8methoxybenzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 6-[2-(4-Methylpiperazin-1-vl)ethyl]benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 6-[2-Morpholin-4-ylethyl]benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 6-(2-Pyrrolidin-1-ylethyl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 6-(2-Piperidin-1-yl)ethyl]benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 4-Bromo-6-[2-morpholin-4-ylethyl]benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 4-Bromo-6-(2-pyrrolidin-1-ylethyl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
- $\label{lem:condition} 4-Bromo-6-[2-(4-methylpiperazin-1-yl)ethyl] benzo[d] isothiazolo[3,2-a] indol-S, S-dioxide;$ 
  - $6\hbox{-}(3\hbox{-}(Piperidin-1\hbox{-}yl)\hbox{-}1\hbox{-}hydroxyprop-1\hbox{-}yl) benzo[d] is othiazolo[3,2\hbox{-}a] indol\hbox{-}S,S-dioxide;$

 $\label{eq:continuous} 6\text{-}(3\text{-}(Piperidin-1\text{-}yl)-1\text{-}hydroxyprop-1\text{-}yl)-8\text{-}methoxybenzo[d]} isothiazolo[3,2\text{-}a]indol-S.S.-dioxide:$ 

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- $\label{lem:condition} 4-Bromo-6-(3-(piperidin-1-yl)-1-hydroxyprop-1-yl) benzo[d] isothiazolo[3,2-a] indol-S, S-dioxide:$
- $\label{lem:condition} 4-Bromo-6-(3-(piperidin-1-yl)-1-hydroxyprop-1-yl)-8-methoxybenzo[d] isothiazolo[3,2-a] indol-S,S-dioxide;$ 
  - 6-(3-(Pyrrolidin-1-yl)-1-hydroxyprop-1-yl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
- 6-(3-(Pyrrolidin-1-yl)-1-hydroxyprop-1-yl)-8-methoxybenzo[d] isothiazolo[3,2-a] indol-8, S-dioxide;
  - 6-(2-(N,N-Diethylamino)-2-methylethyl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
  - 6-(2-(N,N-Dimethylamino-1-hydroxy-1-yl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
- 4-Bromo-6-(2-(N,N-Dimethylamino-1-hydroxy-1-yl)benzo[d]isothiazolo[3,2-a]indol-S,S-dioxide;
- $\label{eq:continuous} 6-(2-(N,N-Dimethylaminoethyl)-2,4-diffuoro-8-M\underline{m}ethoxybenzo[d] is othiazolo[3,2-a] indol-S,S-dioxide;$ 
  - $6\hbox{-}(2\hbox{-}(N,N\hbox{-}Dimethylamino\hbox{-}2\hbox{-}methylethyl) benzo[d] is othiazolo[3,2\hbox{-}a] indol\hbox{-}S,S\hbox{-}dioxide;}$
- $\label{lem:condition} 4- Chloro-6-(2-(N,N-Dimethylaminoethyl)-8-methylbenzo[d] is othiazolo[3,2-a] indol-S, S-dioxide; and$ 
  - 8-(2-(N,N-Dimethylaminoethyl)benzo[d]isothiazolo[3,2-a]benzo(g)indol-S,S-dioxide, or its stereoisomers, its N-oxides, and or its pharmaceutically acceptable salts.
- 3. (Currently amended) A pharmaceutical composition comprising either of a pharmaceutically acceptable carrier, diluent/s, excipient/s or solvents along with a therapeutically

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effective amount of a compound according to Claim 1, its tautomeric forms, its stereoisomers, its geometric forms, its N-oxides, and or its pharmaceutically acceptable salts.

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(Currently Amended) A pharmaceutical composition according to Claim 3, in the form of a
tablet, capsule, powder, lozenges, suppositories, syrup, solution, suspension or injectable,
administered—n. as a single dose or multiple dose units.

## 5-25. (Canceled)

26. (Currently amended) Novel intermediates Intermediates of formula (III) are represented as given below.

$$\begin{array}{c|c}
R_1 & R_9 & H \\
R_{10} & H & H \\
\hline
R_{11} & R_{11} & R_{11} \\
R_{12} & R_8 & R_7 \\
\hline
R_1 & R_9 & R_1
\end{array}$$
(III)

wherein R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, R<sub>8</sub>, R<sub>9</sub>, R<sub>10</sub>, R<sub>11</sub> and R<sub>12</sub> may be the same or different and each independently represent hydrogen, halogen, perhaloalkyl, hydroxy, amino, nitro, cyano, formyl, amidino, guanidino, substituted or unsubstituted groups such as selected from the group consisting of linear or branched (C<sub>1</sub>-C<sub>12</sub>)alkyl, (C<sub>2</sub>-C<sub>12</sub>)alkenyl, (C<sub>2</sub>-C<sub>12</sub>)alkynyl, (C<sub>3</sub>-C<sub>7</sub>)cycloalkenyl, bicycloalkyl, bicycloalkenyl, (C<sub>1</sub>-C<sub>12</sub>)alkoxy, cyclo(C<sub>3</sub>-C<sub>7</sub>)alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heterocyclylalkyl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, heterocyclylalkyloxy, acyl, acyloxy, acylamino, monoalkylamino, dialkylamino, aralkylamino, aralkylamino, alkoxycarbonyl, aryloxycarbonyl, aralkoxycarbonyl, heteroaryloxycarbonyl, heterocyclylalkoxycarbonyl, heteroaryloxycarbonyl, heteroaryloxycarbonyl, aminoalkyl,

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monoalkylaminoalkyl, dialkylaminoalkyl, alkoxyalkyl, aryloxyalkyl, aralkoxyalkyl, alkylthio, alkoxycarbonylamino. aryloxycarbonylamino, aralkyloxycarbonylamino, thioalkyl. aminocarbonylamino. alkylaminocarbonylamino, dialkylaminocarbonylamino, alkylguanidino, dialkylguanidino, hydrazino, and hydroxylamino; or the adjacent groups like R1 and R2 or R2 and R3 or R3 and R4 or R5 and R6 or R6 and R7 or R7 and R8 together with carbon atoms to which they are attached may form a 5, 6, or 7 membered ring, which may further optionally contain one or more double bonds and/or one or more heteroatoms such as selected from the group consisting of the group "Oxygen", "Nitrogen", "Sulfur" or "Selenium" and combinations of double bond and heteroatoms; or Ro and R10 or R11 and R12 together represent double bond attached to "Oxygen" or "Sulfur"; or Ro and R10 or R11 and R12 together with the carbon atoms to which they are attached may form a 3, 4, 5, or 6 membered ring, which may further optionally contain one or more double bonds, and/or one or more heteroatoms such as selected from the group consisting of the group "Oxygen", "Nitrogen", "Sulfur"-or and "Selenium", as above defined;

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"n" is an integer ranging from 1 to 8.

27-29. (Canceled)

- 30. (Previously presented) The compound of claim 1, wherein n is 1 to 4.
- 31. (Currently amended) The novel intermediates of claim 26, wherein n is 1 to 4.